NEWS 7 Apr 22 BIOSIS Gene Names now available in TOXCENTER NEWS 8 Apr 22 Federal Research in Progress (FEDRIP) now available NEWS 9 Jun 03 New e-mail delivery for search results now available NEWS 10 Jun 10 MEDLINE Reload NEWS 11 Jun 10 PCTFULL has been reloaded NEWS 12 Jul 02 FOREGE no longer contains STANDARDS file segment NEWS 13 Jul 22 USAN to be reloaded July 28, 2002; saved answer sets no longer valid NEWS 14 Jul 29 Enhanced polymer searching in REGISTRY NEWS 15 Jul 30 NETFIRST to be removed from STN NEWS 16 Aug 08 CANCERLIT reload NEWS 17 Aug 08 PHARMAMarketLetter(PHARMAML) - new on STN NEWS 18 Aug 08 NTIS has been reloaded and enhanced NEWS 19 Aug 09 JAPIO to be reloaded August 25, 2002 NEWS 20 Aug 19 Aquatic Toxicity Information Retrieval (AQUIRE) now available on STN NEWS 21 Aug 19 IFIPAT, IFICDB, and IFIUDB have been reloaded Aug 19 The MEDLINE file segment of TOXCENTER has been reloaded NEWS 22 NEWS 23 Aug 26 Sequence searching in REGISTRY enhanced NEWS EXPRESS February 1 CURRENT WINDOWS VERSION IS V6.0d, CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP), AND CURRENT DISCOVER FILE IS DATED 05 FEBRUARY 2002 NEWS HOURS STN Operating Hours Plus Help Desk Availability General Internet Information NEWS INTER NEWS LOGIN Welcome Banner and News Items Direct Dial and Telecommunication Network Access to STN NEWS PHONE NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

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FILE 'HOME' ENTERED AT 06:22:05 ON 28 AUG 2002

=> file cacplus
'CACPLUS' IS NOT A VALID FILE NAME
SESSION CONTINUES IN FILE 'HOME'

Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files that are available. If you have requested multiple files, you can specify a corrected file name or you can enter "IGNORE" to continue accessing the remaining file names entered.

=> e
N-[N-[3-(3-methoxy-4-hydroxyphenyl)propyl]-L-.alpha.-aspartyl]-L-phenylalanine
1-Me ester/cn
THIS COMMAND NOT AVAILABLE IN THE CURRENT FILE
The EXPAND command is used to look at the index in a file
which has an index. This file does not have an index.

=> file reg COST IN U.S. DOLLARS

FULL ESTIMATED COST

E12

1

SINCE FILE TOTAL
ENTRY SESSION
0.63 0.63

FILE 'REGISTRY' ENTERED AT 06:23:52 ON 28 AUG 2002 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2002 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6 DICTIONARY FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting ${\tt SmartSELECT}$ searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=> e
N-[N-[3-(3-methoxy-4-hydroxyphenyl)propyl]-L-.alpha.-aspartyl]-L-phenylalanine
1-Me ester/cn
E1 1 N,S-DIVINYL-2-MERCAPTOBENZIMIDAZOLE/CN
E2 1 N,S-TRANS-1,2-DITHIOLANYLIDENETHIAMINE/CN
E3 0 --> N-N-3-(3-METHOXY-4-HYDROXYPHENYL)PROPYL -L-.ALPHA.-ASPARTY
L -L-PHENYLALANINE 1-ME ESTER/CN .

E4 1
N-(((4-(2-((5-CHLORO-2-METHOXYBENZOYL)((4-NITROPHENYL)SULFO

NYL) AMINO) ETHYL) PHENYL) SULFONYL) ((4-NITROPHENYL) SULFONYL) AMI
NO) CARBONYL) -N-CYCLOHEXYL-4-NITROBENZENESULFONAMIDE/CN

N-(((1-(3,4-DIHYDROXY-5-(HYDROXYMETHYL)TETRAHYDRO-2-FURANYL)

-5-IODO-2-OXO-1,2-DIHYDRO-4-PYRIMIDINYL) ((4-NITROPHENYL) SULF

ONYL) AMINO) CARBOTHIOYL) - 4-NITRO-N-(4-NITROPHENYL) BENZENESULF

		ONAPTI DE / CN
E6	1	N-(((2,3-DIMETHOXYBENZYL)OXY)CARBONYL)TRYPTOPHAN/CN
E7	1	N-(((2,6-DICHLOROPHENYL)CARBAMOYL)METHYL)IMINODIACETIC
ACID/		
		CN
E8	1	N-(((2,6-DIISOPROPYLPHENYL)CARBAMOYL)METHYL)IMINODIACETIC
AC		
		ID/CN
E9	1	N-(((2-(DIETHYLAMINO)ETHYL)OXY)CARBONYL)CYCLOPEPTIDE A,
HYDR		
		OCHLORIDE/CN
E10	1	N-(((2-AMINOETHYL)AMINOMETHYL)BENZYLOXY)DIMETHYLSILANE/CN
E11	1	N-(((2-HYDROXYBENZOYL)OXY)METHYL)BENZAMIDE/CN
	-	///2

N-(((3,4-DIMETHOXYPHENYL)METHYLENE)AMINO)PHTHALIMIDE/CN

```
=> e
N-(N-(3-(3-methoxy-4-hydroxyphenyl)propyl)-L-.alpha.-aspartyl)-L-phenylalanine
1-Me ester/cn
                   N-(N-(2-METHYLPHENYL)-N-PHENYLAMINO) PHTHALIMIDE/CN
E1
                   N-(N-(3,4-DICHLOROPHENYL) CARBAMOYL) MORPHOLINE/CN
E2
             1
E3
             0 -->
N-(N-(3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL)-L-.ALPHA.-ASPARTY
                   L)-L-PHENYLALANINE 1-ME ESTER/CN
E4
             1
N-(N-(3-AMINO-3-CARBOXYPROPYL)-3-AMINO-3-CARBOXYPROPYL)AZETI
                   DINE-2-CARBOXYLIC ACID/CN
                   N-(N-(3-CHLOROPHENYL)-N-PHENYLAMINO) PHTHALIMIDE/CN
             1
                   N-(N-(3-PHENOXYBENZOYL)GLYCYL)GLYCINE/CN
E6
             1
E7
                   N-(N-(4-CARBOXYPHENYL)GLYCYL)AMINOACETONITRILE/CN
F.8
N-(N-(5-ETHOXYCARBONYL-4-METHYL-4-PENTENOYL)-L-ALANYL)-L-VAL
                   INE METHYL ESTER/CN
             1
N-(N-(9-.BETA.-D-RIBOFURANOSYLPURIN-6-YL) CARBAMOYL) THREONINE
                   /CN
                   N-(N-(ACRYLAMIDOMETHYL) CARBAMOYLETHYL) PHENOTHIAZINE/CN
E10
             1
             1
                   N-(N-(ADAMANTANE-1-CARBONYL)GLYCYL)GLYCINE/CN
E11
E12
             1
                   N-(N-(BENZYLOXYCARBONYL) GLYCYL) GLYCINE/CN
=>
Uploading 10091500 aspartame derivative.str
L1
```

STRUCTURE UPLOADED

=> d 11

L1 HAS NO ANSWERS

L1 STR

Structure attributes must be viewed using STN Express query preparation.

=> Uploading 10091500 aspartame derivative.str

L2 STRUCTURE UPLOADED => d 12 L2 HAS NO ANSWERS L2 STR

Structure attributes must be viewed using STN Express query preparation.

=> search 12 exact sam

SAMPLE SEARCH INITIATED 06:27:08 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 2 TO ITERATE

100.0% PROCESSED

2 ITERATIONS

0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS:

2 TO 124

PROJECTED ITERATIONS
PROJECTED ANSWERS:

0 TO 12.

- -

0 SEA EXA SAM L2

=> search 12 exact full

FULL SEARCH INITIATED 06:27:15 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 13 TO ITERATE

100.0% PROCESSED

13 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

L4

1 SEA EXA FUL L2

=> d 14

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 329326-75-2 REGISTRY

CN L-Phenylalanine, N-[3-(4-hydroxy-3-methoxyphenyl)propyl]-L-.alpha.aspartyl-, 2-methyl ester (9CI) (CA INDEX NAME)

FS STEREOSEARCH

MF C24 H30 N2 O7

SR CA

LC STN Files: CA, CAPLUS, CASREACT

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

2 REFERENCES IN FILE CA (1967 TO DATE)
2 REFERENCES IN FILE CAPLUS (1967 TO DATE)

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 50.86 51.49

FULL ESTIMATED COST

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=> 14

L5 2 L4

=> d 15 1-2 ti fbib abs ANSWER 1 OF 2 CAPLUS COPYRIGHT 2002 ACS ΤI Process for producing aspartyl dipeptide ester derivatives 2001:833348 CAPLUS AN 135:358168 DN TΙ Process for producing aspartyl dipeptide ester derivatives Kawahara, Shigeru; Nagashima, Kazutaka; Takemoto, Tadashi IN PA Ajinomoto Co., Inc., Japan SO PCT Int. Appl., 25 pp. CODEN: PIXXD2 DT Patent Japanese LA FAN.CNT 1 PATENT NO. KIND DATE APPLICATION NO. DATE WO 2001085761 A1 20011115 WO 2001-JP3479 20010423 PΤ W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG JP 2000-137028 A 20000510 CASREACT 135:358168; MARPAT 135:358168 OS This document discloses a process for conveniently producing on an industrial scale in high yield N-{N-[3-(phenyl)propyl]-L-.alpha.-aspartyl]-L-phenylalanine 1-Me ester derivs., which are expected to be sweeteners, by reductively alkylating aspartame with 3-phenyl-2-propenyl aldehyde derivs. under hydrogen in the presence of a catalyst and a base. RE.CNT 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT L5 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2002 ACS Process for producing and purifying aspartame derivative as sweetener TΙ AN 2001:185780 CAPLUS DN 134:223039 Process for producing and purifying aspartame derivative as sweetener ΤI Amino, Yusuke; Yuzawa, Kazuko; Takemoto, Tadashi IN Ajinomoto Co., Inc., Japan PA SO PCT Int. Appl., 39 pp. CODEN: PIXXD2 DT Patent Japanese LA FAN.CNT 1 PATENT NO. KIND DATE APPLICATION NO. DATE ______ WO 2001018034 A1 20010315 WO 2000-JP5665 20000823 PT W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR,

HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN,

YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM

RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

JP 1999-253498 A 19990907

OS CASREACT 134:223039

AB This document discloses the following: a method for industrially producing

N-[N-[3-(3-methoxy-4-hydroxyphenyl)propyl]-L-.alpha.-aspartyl]-L-phenylalanine l-Me ester which is useful as a sweetener, in particular, a process for producing the target compd. in a high yield by the reductive alkylation reaction of aspartame with 3-(3-methoxy-4-hydroxyphenyl)propional dehyde or its deriv.; a method of effectively purifying the target compd. contaminated with impurities invading thereinto at various prodn. stages (involving methods other than the above-described reductive alkylation), more particularly, a method of sepg. the target compd. in the form of highly pure crystals; the crystals;

sweeteners contg. the same; and utilization thereof in various products which are to be sweetened.

RE.CNT 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> logoff hold SINCE FILE TOTAL COST IN U.S. DOLLARS ENTRY SESSION 5.37 56.86 FULL ESTIMATED COST SINCE FILE DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) TOTAL ENTRY SESSION -1.24-1.24CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 60 MINUTES STN INTERNATIONAL SESSION SUSPENDED AT 06:28:46 ON 28 AUG 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1623paz

PASSWORD:

* * * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * * SESSION RESUMED IN FILE 'CAPLUS' AT 07:09:05 ON 28 AUG 2002 FILE 'CAPLUS' ENTERED AT 07:09:05 ON 28 AUG 2002 COPYRIGHT (C) 2002 AMERICAN CHEMICAL SOCIETY (ACS)

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	5.37	56.86
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-1.24	-1.24
=> file reg		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL

SESSION ENTRY 56,86 5.37 FULL ESTIMATED COST TOTAL DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE ENTRY SESSION -1.24 -1.24CA SUBSCRIBER PRICE FILE 'REGISTRY' ENTERED AT 07:09:10 ON 28 AUG 2002 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2002 American Chemical Society (ACS) STRUCTURE FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6 DICTIONARY FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6 TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002 Please note that search-term pricing does apply when conducting SmartSELECT searches. Crossover limits have been increased. See HELP CROSSOVER for details. Calculated physical property data is now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf => e 3-(4-hydroxy-3-methoxyphenyl) propionaldehyde/cn 3-(4-HYDROXY-3-METHOXYBENZYL)-5-HYDROXY-6,7-DIMETHOXYCHROMAN -4-ONE/CN 3-(4-HYDROXY-3-METHOXYBENZYL)-5-HYDROXY-7-METHOXYCHROMAN-4-0 NE/CN 0 --> 3-(4-HYDROXY-3-METHOXYPHENYL) PROPIONALDEHYDE/CN 3-(4-HYDROXY-3-METHOXYPHENYL)-1,2-PROPANEDIOL/CN E4 1 3-(4-HYDROXY-3-METHOXYPHENYL)-1,2-PROPANEDIOL TRIACETATE/CN E5 1 3-(4-HYDROXY-3-METHOXYPHENYL)-1-PHENYL-1-PROPENE/CN F.6 1 3-(4-HYDROXY-3-METHOXYPHENYL)-1-PROPANOL/CN E7 1 3-(4-HYDROXY-3-METHOXYPHENYL)-1-PROPENE/CN 1 E.S FQ 1 3-(4-HYDROXY-3-METHOXYPHENYL)-2-(2-METHOXY-4-METHYLPHENOXY)-2-PROPEN-1-OL/CN 1 E10 3-(4-HYDROXY-3-METHOXYPHENYL)-2-(2-METHOXY-4-PROPYLPHENOXY)-1-PROPANOL/CN 1 3-(4-HYDROXY-3-METHOXYPHENYL)-2-(2-METHOXY-5-METHYLPHENOXY)-2-PROPEN-1-OL/CN E12 3-(4-HYDROXY-3-METHOXYPHENYL)-2-(3,4-DIMETHOXYPHENYL)-2-PROP ENESULFONIC ACID/CN => e 3-(4-hydroxy-3-methoxyphenyl)-propionaldehyde/cn 3-(4-HYDROXY-3-METHOXYPHENYL)-2-PROPENOIC ACID/CN E.1 3-(4-HYDROXY-3-METHOXYPHENYL)-5-PHENYL-1,2-DITHIOLIUM E2 7 PERCHL

0 --> 3-(4-HYDROXY-3-METHOXYPHENYL)-PROPIONALDEHYDE/CN

ORATE/CN

E3

E4	1	3-(4-HYDROXY-3-METHOXYPHENYL)ACROLEIN/CN
E5	1	3-(4-HYDROXY-3-METHOXYPHENYL)ACRYLIC ACID/CN
E6	1	3-(4-HYDROXY-3-METHOXYPHENYL)ALANINE/CN
E7	1	3-(4-HYDROXY-3-METHOXYPHENYL)ALLYL ALCOHOL/CN
E8	1	3-(4-HYDROXY-3-METHOXYPHENYL)LACTIC ACID/CN
E9	1	3-(4-HYDROXY-3-METHOXYPHENYL)PROPA-1,2-DIENE/CN
E10	1	3-(4-HYDROXY-3-METHOXYPHENYL)PROPANOL/CN
E11	1	3-(4-HYDROXY-3-METHOXYPHENYL) PROPIONIC ACID/CN
E12	1	3-(4-HYDROXY-3-METHOXYSTYRYL)-2,5-DIMETHYLISOXAZOLIUM
IODIDE		
		/CN

=> logoff hold

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 1.90 58.76

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION 0.00 -1.24CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 60 MINUTES STN INTERNATIONAL SESSION SUSPENDED AT 07:11:53 ON 28 AUG 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1623paz

PASSWORD:

* * * * * * RECONNECTED TO STN INTERNATIONAL * * * * * SESSION RESUMED IN FILE 'REGISTRY' AT 07:17:02 ON 28 AUG 2002 FILE 'REGISTRY' ENTERED AT 07:17:02 ON 28 AUG 2002 COPYRIGHT (C) 2002 American Chemical Society (ACS)

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	1.90	58.76
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	0.00	-1.24

Uploading 10091500 aldehyde.str

L6 STRUCTURE UPLOADED

=> d 16

L6 HAS NO ANSWERS L6 STR

Structure attributes must be viewed using STN Express query preparation.

=> search 16 exact full FULL SEARCH INITIATED 07:18:00 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 124 TO ITERATE

100.0% PROCESSED 124 ITERATIONS 1 ANSWERS SEARCH TIME: 00.00.01

L7 1 SEA EXA FUL L6

=> d 17

L7 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 80638-48-8 REGISTRY

CN Benzenepropanal, 4-hydroxy-3-methoxy- (9CI) (CA INDEX NAME)

OTHER NAMES:

CN 3-(4'-Hydroxy-3'-methoxyphenyl)-1-propanal

CN 4-Propanal-2-methoxyphenol

FS 3D CONCORD MF C10 H12 O3

LC STN Files: CA, CAPLUS, CASREACT, CHEMLIST, TOXCENTER

Other Sources: DSL**, EINECS**

(**Enter CHEMLIST File for up-to-date regulatory information)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 11 REFERENCES IN FILE CA (1967 TO DATE)
- 11 REFERENCES IN FILE CAPLUS (1967 TO DATE)

=> file caplus TOTAL COST IN U.S. DOLLARS SINCE FILE ENTRY SESSION FULL ESTIMATED COST 51.24 108.10 SINCE FILE TOTAL DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -1.24

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=> 17 L8 11 L7

=> save temp 18 sweetaldeh/a ANSWER SET L8 HAS BEEN SAVED AS 'SWEETALDEH/A'

=>
COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)
CA SUBSCRIBER PRICE
SINCE FILE TOTAL
ENTRY SESSION
0.00 -1.24

FILE 'STNGUIDE' ENTERED AT 07:19:02 ON 28 AUG 2002
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FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Aug 23, 2002 (20020823/UP).

=>			
NAME	CREATED	NOTES/TITLE	
ADIACRYST/A ALKYLATIN/L			
ASPARTAME/A A408SRCH/L	TEMP	2346 ANSWERS IN FILE CAPLUS	
CONIFERAL/A GLYPHSRCH/L	TEMP	388 ANSWERS IN FILE CAPLUS	
INDIUMCL3/A	30 MAY 2001	1 ANSWER IN FILE REGISTRY 72 ANSWERS IN FILE CAPLUS	
NASTATINS/A NEOTAMECRYST/A	24 APR 2001	59 ANSWERS IN FILE CAPLUS	
POHBENZALDEH/A	10 JUL 2001	196 ANSWERS IN FILE REGISTRY 5519 ANSWERS IN FILE CAPLUS	
PROCTYLCMPD/A PROCTYLSRCH/L	TEMP		
STILLEAPP/L	07 JAN 2002 TEMP TEMP	17 L-NUMBERS 6 L-NUMBERS 11 ANSWERS IN FILE CAPLUS	
TWOAMINOPOLY/Q	16 APR 2001	UPLOADED STRUCTURE	

=>
NO SAVED SDI REQUESTS

SINCE FILE TOTAL COST IN U.S. DOLLARS ENTRY SESSION 0.06 108.56 FULL ESTIMATED COST DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL SESSION ENTRY CA SUBSCRIBER PRICE 0.00 -1.24

FILE 'CAPLUS' ENTERED AT 07:19:47 ON 28 AUG 2002 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
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=> L9 (L10

1)SEA FILE=REGISTRY ABB=ON PLU=ON ASPARTAME/CN

2346 SEA FILE=CAPLUS ABB=ON PLU=ON L9

=>

L11 (1) SEA FILE=REGISTRY ABB=ON PLU=ON CONIFERALDEHYDE/CN

L12 388 SEA FILE=CAPLUS ABB=ON PLU=ON L11

=> file caplus

COST IN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST 0.40 108.96

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SI

SINCE FILE TOTAL ENTRY SESSION

CA SUBSCRIBER PRICE

0.00 -1.24

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=> d his

(FILE 'HOME' ENTERED AT 06:22:05 ON 28 AUG 2002)

FILE 'REGISTRY' ENTERED AT 06:23:52 ON 28 AUG 2002

N-[N-[3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL]-L-.ALPHA.-ASPARTYL

N-(N-(3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL)-L-.ALPHA.-ASPARTYL
L1 STRUCTURE UPLOADED

```
STRUCTURE UPLOADED
L2
L3
              0 SEARCH L2 EXACT SAM
L4
              1 SEARCH L2 EXACT FULL
     FILE 'CAPLUS' ENTERED AT 06:27:35 ON 28 AUG 2002
L5
              2 L4
     FILE 'REGISTRY' ENTERED AT 07:09:10 ON 28 AUG 2002
                E 3-(4-HYDROXY-3-METHOXYPHENYL) PROPIONALDEHYDE/CN
                E 3-(4-HYDROXY-3-METHOXYPHENYL)-PROPIONALDEHYDE/CN
L6
                STRUCTURE UPLOADED
L7
              1 SEARCH L6 EXACT FULL
     FILE 'CAPLUS' ENTERED AT 07:18:17 ON 28 AUG 2002
L8
             11 L7
                SAVE TEMP L8 SWEETALDEH/A
     FILE 'STNGUIDE' ENTERED AT 07:19:02 ON 28 AUG 2002
     FILE 'CAPLUS' ENTERED AT 07:19:47 ON 28 AUG 2002
               ACT ASPARTAME/A
L9 (
              1) SEA FILE=REGISTRY ABB=ON PLU=ON ASPARTAME/CN
L10
           2346 SEA FILE=CAPLUS ABB=ON PLU=ON L9
               ACT CONIFERAL/A
               -----
L11 (
              1) SEA FILE=REGISTRY ABB=ON PLU=ON CONIFERALDEHYDE/CN
L12
            388 SEA FILE=CAPLUS ABB=ON PLU=ON L11
               _____
     FILE 'CAPLUS' ENTERED AT 07:20:01 ON 28 AUG 2002
=> 18 and 110
L13
            1 L8 AND L10
=> d 113 ti fbib abs
L13 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2002 ACS
    Process for producing and purifying aspartame derivative as sweetener
TΙ
AN
     2001:185780 CAPLUS
    134:223039
DN
TI
    Process for producing and purifying aspartame derivative as sweetener
    Amino, Yusuke; Yuzawa, Kazuko; Takemoto, Tadashi
IN
    Ajinomoto Co., Inc., Japan
PΑ
SO
     PCT Int. Appl., 39 pp.
    CODEN: PIXXD2
DT
    Patent
LΑ
    Japanese
FAN.CNT 1
                 KIND DATE
    PATENT NO.
                                         APPLICATION NO. DATE
                     ____
    WO 2001018034 A1 20010315
                                         WO 2000-JP5665 20000823
PΤ
        W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
             CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR,
             HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT,
             LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU,
             SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
```

RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

JP 1999-253498 A 19990907

OS CASREACT 134:223039

AB This document discloses the following: a method for industrially producing

N-[N-[3-(3-methoxy-4-hydroxyphenyl)propyl]-L-.alpha.-aspartyl]-Lphenylalanine 1-Me ester which is useful as a sweetener, in particular, a
process for producing the target compd. in a high yield by the reductive
alkylation reaction of aspartame with 3-(3-methoxy-4hydroxyphenyl)propionaldehyde or its deriv.; a method of effectively
purifying the target compd. contaminated with impurities invading
thereinto at various prodn. stages (involving methods other than the
above-described reductive alkylation), more particularly, a method of
sepg. the target compd. in the form of highly pure crystals; the
crystals;

sweeteners contg. the same; and utilization thereof in various products which are to be sweetened.

RE.CNT 11 THERE ARE 11 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> 18 and 112 L14 6 L8 AND L12

=> d 114 1-6 ti

1.14 ANSWER 1 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Process for producing and purifying aspartame derivative as sweetener

L14 ANSWER 2 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Retention of lignin in seagrasses: angiosperms that returned to the sea

L14 ANSWER 3 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Chemical composition of alkaline pulping products of lignin model compounds

L14 ANSWER 4 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Thermolytic decomposition of coniferyl alcohol

L14 ANSWER 5 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Ultrafiltration and pyrolysis gas chromatography mass spectrometry of chlorolignins in pulp mill effluent

L14 ANSWER 6 OF 6 CAPLUS COPYRIGHT 2002 ACS

TI Topical cosmetics containing 1,7-diphenyl-4-hepten-3-one for skin disorder

treatment

=> logoff hold SINCE FILE TOTAL COST IN U.S. DOLLARS ENTRY SESSION 122.06 FULL ESTIMATED COST 13.10 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL SESSION ENTRY -0.62 -1.86 CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 60 MINUTES STN INTERNATIONAL SESSION SUSPENDED AT 07:33:50 ON 28 AUG 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1623paz

PASSWORD:

* * * * * * RECONNECTED TO STN INTERNATIONAL * * * * * * SESSION RESUMED IN FILE 'CAPLUS' AT 07:51:23 ON 28 AUG 2002 FILE 'CAPLUS' ENTERED AT 07:51:23 ON 28 AUG 2002 COPYRIGHT (C) 2002 AMERICAN CHEMICAL SOCIETY (ACS)

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	, 13.10	122.06
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-0.62	-1.86
=> file reg		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	13.89	122.85
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-0.62	-1.86

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STRUCTURE FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6 DICTIONARY FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=> e Benzenepropanal, 4-hydroxy-3-methoxy-, cn

E1 106272 BENZENEPROPAN/BI

E2 2540 BENZENEPROPANAL/BI

E3 0 --> BENZENEPROPANAL, 4-HYDROXY-3-METHOXY-, CN/BI

E4	60	BENZENEPROPANALATO/BI		
E5	1	BENZENEPROPANALDEHYDE/BI		
E6	538	BENZENEPROPANAMID/BI		
E7	536	BENZENEPROPANAMIDATO/BI		
E8	24838	BENZENEPROPANAMIDE/BI		
E9	1	BENZENEPROPANAMIDO/BI		
E10		BENZENEPROPANAMIN/BI		
E11		BENZENEPROPANAMINATO/BI		
E12	5004	BENZENEPROPANAMINE/BI		
=> e 4	-hydroxy-3-	methoxy-Benzenepropanal cn		
E1	î	4-AMINO-2,6-PYRIMIDINODITHI	OCARBAMATE/BI	
E2	1			
4-HYDR	OXY-2-METHY	L-N-2-PYRIDINYL-2H-1,2-BENZOT	HIAZINE-3-CARB	
		OXAMIDE 1,1-DIOXIDE-NN/BI		
E3	0	> 4-HYDROXY-3-METHOXY-BENZENE	PROPANAL CN/BI	
E4	3	4-ISOCYANATOBENZENE/BI		
E5	1	4-METHOXYALPHAMETHYLBEN	IZENEACETATE/BI	
E6	2	4-METHYL-2-OXO-2H-1-BENZOPY 4-METHYLBENZENESULFONAMIDAT	RAN-7-YL/BI	
E7	1	4-METHYLBENZENESULFONAMIDAT	O/BI	
E8	1	4-METHYLBENZENESULFONATO P	KAPPA.O/BI	
E9	1	4-METHYLPHENYL		
6-DIAZ	D-5,6-DIHYD	RO-5-OXO-1-NAPHTHALENESULFON		
		ATE/BI		
E10	1	4-MORPHOLINO-1,3,5-TRIAZINE	E-2,6-DIYL/BI	
E11		4-NITROPHENYL/BI		
E12	1	4-0BETAD-GALACTOPYRANOS	SYL-D-GLUCONATO-	01,02,03/BI
=> fil	e caplus			
COST I	U.S. DOLL	ARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST 1.14 123.99			123.99	
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION				
CA SUB	SCRIBER PRI	CE	0.00	-1.86

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=> d his

(FILE 'HOME' ENTERED AT 06:22:05 ON 28 AUG 2002)

FILE 'REGISTRY' ENTERED AT 06:23:52 ON 28 AUG 2002

N-[N-[3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL]-L-.ALPHA.-ASPARTYL Ε

N-(N-(3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL)-L-.ALPHA.-ASPARTYL

L1 STRUCTURE UPLOADED

L2 STRUCTURE UPLOADED

L3 0 SEARCH L2 EXACT SAM

1 SEARCH L2 EXACT FULL L4

FILE 'CAPLUS' ENTERED AT 06:27:35 ON 28 AUG 2002

2 L4 L5

FILE 'REGISTRY' ENTERED AT 07:09:10 ON 28 AUG 2002

E 3-(4-HYDROXY-3-METHOXYPHENYL) PROPIONALDEHYDE/CN

E 3-(4-HYDROXY-3-METHOXYPHENYL)-PROPIONALDEHYDE/CN

L6 STRUCTURE UPLOADED

1 SEARCH L6 EXACT FULL T.7

FILE 'CAPLUS' ENTERED AT 07:18:17 ON 28 AUG 2002

T.R 11 L7

SAVE TEMP L8 SWEETALDEH/A

FILE 'STNGUIDE' ENTERED AT 07:19:02 ON 28 AUG 2002

FILE 'CAPLUS' ENTERED AT 07:19:47 ON 28 AUG 2002

ACT ASPARTAME/A

1) SEA FILE=REGISTRY ABB=ON PLU=ON ASPARTAME/CN L9 (

L10 2346 SEA FILE=CAPLUS ABB=ON PLU=ON L9

ACT CONIFERAL/A

1) SEA FILE=REGISTRY ABB=ON PLU=ON CONIFERALDEHYDE/CN L11 (

388 SEA FILE=CAPLUS ABB=ON PLU=ON L11 L12

FILE 'CAPLUS' ENTERED AT 07:20:01 ON 28 AUG 2002

1 L8 AND L10 L13

L14 6 L8 AND L12

FILE 'REGISTRY' ENTERED AT 07:52:26 ON 28 AUG 2002

E BENZENEPROPANAL, 4-HYDROXY-3-METHOXY-, CN

E 4-HYDROXY-3-METHOXY-BENZENEPROPANAL CN

FILE 'CAPLUS' ENTERED AT 07:54:07 ON 28 AUG 2002

=> reg

915 REG

41 REGS

L15 952 REG

(REG OR REGS)

=> file reg

SINCE FILE TOTAL ENTRY SESSION COST IN U.S. DOLLARS 2.02 126.01 FULL ESTIMATED COST DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -1.86

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26 AUG 2002 HIGHEST RN 444986-65-6 STRUCTURE FILE UPDATES: DICTIONARY FILE UPDATES: 26 AUG 2002 HIGHEST RN 444986-65-6

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=> e conife	=> e coniferaldehyde/cn		
E1	1 CONIDIOSPORE SURFACE PROTEIN (TRICHODERMA HARZIANUM STRAIN		
A			
	TCC 32173 GENE CMP1)/CN		
E2	1 CONIFEGEROL/CN		
E3	1> CONIFERALDEHYDE/CN		
E4	1 CONIFERALDEHYDE METHOXYMETHYL ETHER/CN		
E5	1 CONIFERALDEHYDE METHYL ETHER/CN		
E6	1 CONIFERAN/CN		
E7	1 CONIFERIC ACID/CN		
E8	1 CONIFERIN/CN		
E9	1 CONIFERIN .BETAGLUCOSIDASE/CN		
E10	1 CONIFERIN PENTAACETATE/CN		
E11	1 CONIFERIN, DIHYDRATE/CN		
E12	1 CONIFERIN-HYDROLYZING .BETAGLUCOSIDASE/CN		
=> e3			
1.16	1 CONTFERALDEHYDE/CN		

=> d 116

L16 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS 458-36-6 REGISTRY 2-Propenal, 3-(4-hydroxy-3-methoxyphenyl)- (9CI) (CA INDEX NAME) OTHER CA INDEX NAMES: CN Cinnamaldehyde, 4-hydroxy-3-methoxy- (6CI, 8CI) OTHER NAMES:

CN 2-Methoxy-4-(3-oxo-1-propenyl)phenol CN 3-(4-Hydroxy-3-methoxyphenyl)acrolein CN 3-Methoxy-4-hydroxycinnamaldehyde CN 4-Hydroxy-3-methoxycinnamaldehyde CN Coniferaldehyde CN Coniferyl aldehyde CN Ferulaldehyde CN Ferulyl aldehyde CN p-Coniferaldehyde FS 3D CONCORD MF C10 H10 O3 CI COM AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS, LC STN Files: BIOTECHNO, CA, CAOLD, CAPLUS, CASREACT, CHEMCATS, CHEMLIST, CSCHEM, DDFU, DRUGU, EMBASE, HODOC*, MEDLINE, NAPRALERT, NIOSHTIC, PIRA, SPECINFO, TOXCENTER, USPATFULL (*File contains numerically searchable property data) Other Sources: EINECS** (**Enter CHEMLIST File for up-to-date regulatory information)

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

388 REFERENCES IN FILE CA (1967 TO DATE)
3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
388 REFERENCES IN FILE CAPLUS (1967 TO DATE)
28 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

=> file capls
'CAPLS' IS NOT A VALID FILE NAME
SESSION CONTINUES IN FILE 'REGISTRY'
Enter "HELP FILE NAMES" at an arrow;

Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files that are available. If you have requested multiple files, you can specify a corrected file name or you can enter "IGNORE" to continue accessing the remaining file names entered.

=> file caplus COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 5.96 131.97 DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL ENTRY SESSION CA SUBSCRIBER PRICE 0.00 -1.86

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N-(N-(3-(3-METHOXY-4-HYDROXYPHENYL) PROPYL)-L-.ALPHA.-ASPARTYL

1 STRUCTURE UPLOADED
2 STRUCTURE UPLOADED

L2 STRUCTURE UPLOADED
L3 0 SEARCH L2 EXACT SAM

L4 1 SEARCH L2 EXACT FULL

FILE 'CAPLUS' ENTERED AT 06:27:35 ON 28 AUG 2002

L5 2 L4

FILE 'REGISTRY' ENTERED AT 07:09:10 ON 28 AUG 2002

E 3-(4-HYDROXY-3-METHOXYPHENYL) PROPIONALDEHYDE/CN E 3-(4-HYDROXY-3-METHOXYPHENYL)-PROPIONALDEHYDE/CN

L6 STRUCTURE UPLOADED

L7 1 SEARCH L6 EXACT FULL

FILE 'CAPLUS' ENTERED AT 07:18:17 ON 28 AUG 2002

L8 11 L7

SAVE TEMP L8 SWEETALDEH/A

FILE 'STNGUIDE' ENTERED AT 07:19:02 ON 28 AUG 2002

FILE 'CAPLUS' ENTERED AT 07:19:47 ON 28 AUG 2002 ACT ASPARTAME/A

L9 (1) SEA FILE=REGISTRY ABB=ON PLU=ON ASPARTAME/CN

L10 · 2346 SEA FILE=CAPLUS ABB=ON PLU=ON L9

ACT CONIFERAL/A

L11 (1) SEA FILE=REGISTRY ABB=ON PLU=ON CONIFERALDEHYDE/CN

L12 388 SEA FILE=CAPLUS ABB=ON PLU=ON L11

FILE 'CAPLUS' ENTERED AT 07:20:01 ON 28 AUG 2002

L13 1 L8 AND L10 L14 6 L8 AND L12

FILE 'REGISTRY' ENTERED AT 07:52:26 ON 28 AUG 2002

E BENZENEPROPANAL, 4-HYDROXY-3-METHOXY-,CN

E 4-HYDROXY-3-METHOXY-BENZENEPROPANAL CN

FILE 'CAPLUS' ENTERED AT 07:54:07 ON 28 AUG 2002

L15 952 REG

FILE 'REGISTRY' ENTERED AT 07:54:37 ON 28 AUG 2002 E CONIFERALDEHYDE/CN

L16 1 E3

FILE 'CAPLUS' ENTERED AT 07:55:19 ON 28 AUG 2002

=> rhodium

52919 RHODIUM 30 RHODIUMS

L17 52920 RHODIUM

(RHODIUM OR RHODIUMS)

=> 112 and 117

L18 1 L12 AND L17

=> d 118 ti

L18 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2002 ACS

TI Process for producing and purifying aspartame derivative as sweetener

=> logoff hold

COST ÎN U.S. DOLLARS
SINCE FILE TOTAL
ENTRY SESSION
FULL ESTIMATED COST 2.70 134.67

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE
ENTRY
SESSION
CA SUBSCRIBER PRICE

0.00
-1.86

SESSION WILL BE HELD FOR 60 MINUTES
STN INTERNATIONAL SESSION SUSPENDED AT 07:56:45 ON 28 AUG 2002

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:ssspta1623paz